ethyleneoxy units; or L_1 is —NRC(=O)(CH₂)_n—, —NRC(=O)(CH₂)_nC(=O)NH—, or —NR $(CH_2)_nC(=O)NH(CH_2)_n$, L_2 is $-(CH_2)_nO$ —, and L_3 is $-(CH_2)_n$ —, where each n is an integer from 1 to 12;

X is an amino acid, a polypentide, a nucleoside, a nucleotide, a polynucleotide, or a protected form thereof; or X is an acid-labile protecting group;

Z is selected from H, CO₂H, OH, NH₂, NHR, NR₂, SH, OP(NR₁R₂)(OR₃), an ester, a cleavable linker, a solid support, a reactive linking group, and a label selected from a fluorescent dye, a hybridization-stabilizing moiety, a chemiluminescent dye, and an affinity ligand, where R1 and R_2 are C_1 – C_{12} alkyl; C_5 – C_{14} aryl; or cycloalkyl containing up to 10 carbon atoms, or when R_1 and R₂ are taken together with the phosphoramidite nitrogen atom, R₁ and R₂ are C₄-C₁₁ alkyldiyl, and R3 is a phosphite ester protecting group; and

Q is selected from the diazo structures:

wherein Ar is C₅-C₁₄ aryl; one of the aryl carbons of the diazo structures is the site of attachment to L1; at least one aryl carbon of each diazo structure is substituted with an electronwithdrawing group and at least one aryl carbon of each diazo structure is substituted with an electron-donating group.

REMARKS

Reconsideration of the application is respectfully requested. By this Preliminary Amendment, non-elected claims 26-75 have been canceled, and claim 1 has been amended. Claims 1-25 are pending.